

# Package ‘gamreg’

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**Type** Package

**Title** Robust and Sparse Regression via Gamma-Divergence

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**Description** Robust regression via gamma-divergence with L1, elastic net and ridge.

**License** GPL (>= 2)

**LazyData** TRUE

**Suggests** mvtnorm

**Imports** glmnet, robustHD,foreach,doParallel

**LinkingTo** Rcpp, RcppArmadillo

**RoxygenNote** 5.0.1

**NeedsCompilation** yes

**Repository** CRAN

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cv.gam

*Robust Cross-Validation***Description**

Compute Robust Cross-Validation for selecting best model.

**Usage**

```
cv.gam(X, Y, init.mode = c("sLTS", "RLARS", "RANSAC"),
       lambda.mode = "lambda0", lmax = 1, lmin = 0.05, nlambda = 50,
       fold = 10, ncores = 1, gam = 0.1, gam0 = 0.5, intercept = "TRUE",
       alpha = 1, ini.subsamp = 0.2, ini.cand = 1000, alpha.LTS = 0.75,
       nlambda.LTS = 40)
```

**Arguments**

X	Predictor variables Matrix.
Y	Response variables Matrix.
init.mode	"sLTS": a initial point is the estimate of sparse least trimmed squares. "RLARS": a initial point is the estimate of Robust LARS. "RANSAC": a initial point is the estimate of RANSAC algorithm.
lambda.mode	"lambda0": Robust Cross-Validation uses grids on range $[0.05\lambda_0, \lambda_0]$ with log scale, where $\lambda_0$ is an estimator of sparse tuning parameter which would shrink regression coefficients to zero.
lmax	When lambda.mode is not lambda0, upper bound of range of grids is lmax.
lmin	When lambda.mode is not lambda0, lower bound of range of grids is lmin.
nlambda	The number of grids for Robust Cross-Validation.
fold	the number of folds for K-fold Robust Cross-Validation. If fold equals to sample size, Robust Cross-Validation is leave-one-out method.
ncores	positive integer giving the number of processor cores to be used for parallel computing (the default is 1 for no parallelization).
gam	Robust tuning parameter of gamma-divergence for regression.
gam0	tuning parameter of Robust Cross-Validation.
intercept	Should intercept be fitted TRUE or set to zero FALSE
alpha	The elasticnet mixing parameter, with $0 \leq \alpha \leq 1$ . alpha=1 is the lasso penalty, and alpha=0 the ridge penalty.
ini.subsamp	The fraction of subsamples in "RANSAC".
ini.cand	The number of candidates for estimating itnial points in "RANSAC".
alpha.LTS	The fraction of subsamples for trimmed squares in "sLTS".
nlambda.LTS	The number of grids for sparse tuning parameter in "sLTS".

## Details

If the "RANSAC" is used as the initial point, the parameter `ini.subsamp` and `ini.cand` can be determined carefully. The smaller `ini.subsamp` is, the more robust initial point is. However, less efficiency.

## Value

<code>lambda</code>	A numeric vector giving the values of the penalty parameter.
<code>fit</code>	All results at each lambda.
<code>Rocv</code>	The result of best model by Robust Cross-Validation.

## Author(s)

Takayuki Kawashima

## References

Kawashima, T. and Fujisawa, H. (2017). *Robust and Sparse Regression via gamma-divergence*, *Entropy*, 19(11).

Fujisawa, H. and Eguchi, S. (2008). *Robust parameter estimation with a small bias against heavy contamination*, *Journal of Multivariate Analysis*, 99(9), 2053-2081.

## Examples

```
## generate data
library(mvtnorm)
n <- 30                # number of observations
p <- 10               # number of explanatory variables

epsilon <- 0.1        # contamination ratio

beta0 <- 0.0          # intercept
beta <- c(numeric(p)) # regression coefficients
beta[1] <- 1
beta[2] <- 2
beta[3] <- 3
beta[4] <- 4

Sigma <- 0.2^t(sapply(1:p, function(i, j) abs(i-j), 1:p))
X <- rmvnorm(n, sigma=Sigma) # explanatory variables
e <- rnorm(n) # error terms

i <- 1:ceiling(epsilon*n) # index of outliers
e[i] <- e[i] + 20        # vertical outliers
Y <- beta0*(numeric(n)+1) + X*%beta

res <- cv.gam(X,Y,nlambda = 5, nlambda.LTS=20 ,init.mode="sLTS")
```

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